

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101216\
 Data File : BG024312.D
 Acq On : 12 Oct 2016 22:02
 Operator : UM/SJ
 Sample : H5164-10
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BDNN2

Quant Time: Oct 13 01:34:47 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG0100716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 13 01:34:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	148492	20.00	ng/ul	0.00
18) Naphthalene-d8	11.12	136	662347	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.91	164	564106	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.64	188	1167821	20.00	ng/ul	0.00
75) Chrysene-d12	21.94	240	1296233	20.00	ng/ul	0.00
83) Perylene-d12	25.39	264	1303419	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.66	96	5075	1.62	ng/uL	0.01
5) Phenol-d5	7.42	99	92514	6.68	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.60	67	235133	25.11	ng/ul	0.00
9) 2-Chlorophenol-d4	7.81	132	204357	21.42	ng/ul	0.00
13) 4-Methylphenol-d8	8.97	113	164484	14.69	ng/ul	0.00
19) Nitrobenzene-d5	9.46	128	136796	29.10	ng/ul	0.00
22) 2-Nitrophenol-d4	10.18	143	152486	29.00	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.72	165	276334	24.13	ng/ul	0.00
29) 4-Chloroaniline-d4	11.25	131	4377	0.36	ng/ul	0.00
43) Dimethylphthalate-d6	14.30	166	1068266	27.12	ng/ul	0.00
46) Acenaphthylene-d8	14.60	160	1248776	27.68	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	40636	5.82	ng/ul	0.00
57) Fluorene-d10	15.89	176	1018298	27.34	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.99	200	196711	28.03	ng/ul	0.00
70) Anthracene-d10	17.74	188	1514137	28.70	ng/ul	0.00
76) Pyrene-d10	20.01	212	1673811	27.67	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.15	264	1689835	28.82	ng/ul	0.00

Target Compounds Ovalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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